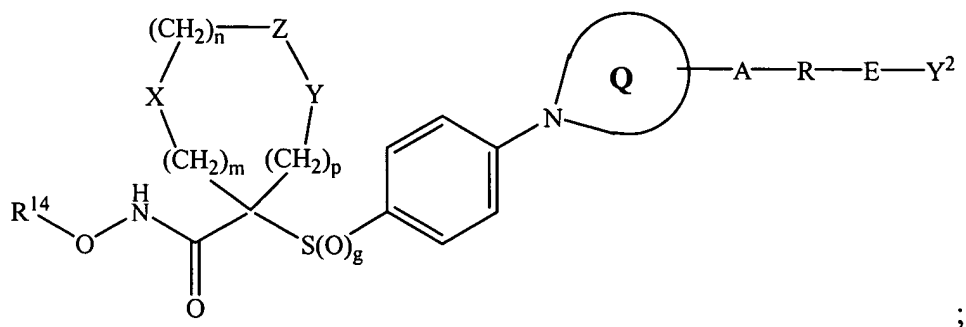


Amended Claims

Claims 1-66 (canceled).

Please add the following new claims:

67. (new) A compound or a pharmaceutically acceptable salt thereof, wherein:
 the compound corresponds in structure to the following formula:



R^{14} is selected from the group consisting of hydrogen and $C(W)R^{25}$;

W is selected from the group consisting of O and S;

R^{25} is selected from the group consisting of C_1 - C_6 -alkyl, aryl, C_1 - C_6 -alkoxy, heteroaryl- C_1 - C_6 -alkyl, C_3 - C_8 -cycloalkyl- C_1 - C_6 -alkyl, aryloxy, aryl- C_1 - C_6 -alkoxy, aryl- C_1 - C_6 -alkyl, heteroaryl, and amino C_1 - C_6 -alkyl, wherein:

the amino C_1 - C_6 -alkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C_1 - C_6 -alkyl, aryl, aryl- C_1 - C_6 -alkyl, C_3 - C_8 -cycloalkyl- C_1 - C_6 -alkyl, aryl- C_1 - C_6 -alkoxycarbonyl, C_1 - C_6 -alkoxycarbonyl, and C_1 - C_6 -alkanoyl, or

the amino C_1 - C_6 -alkyl nitrogen and two substituents attached thereto form a 5- to 8-membered heterocyclo or heteroaryl ring;

g is zero, 1, or 2;

m is zero, 1, or 2;

n is zero, 1, or 2;

p is zero, 1, or 2;

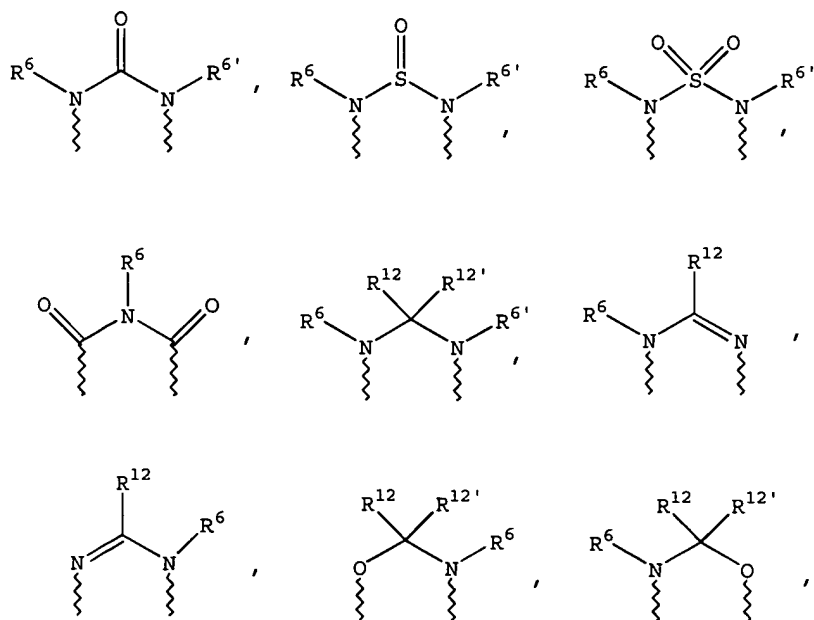
the sum of $m + n + p = 1, 2, 3$, or 4;

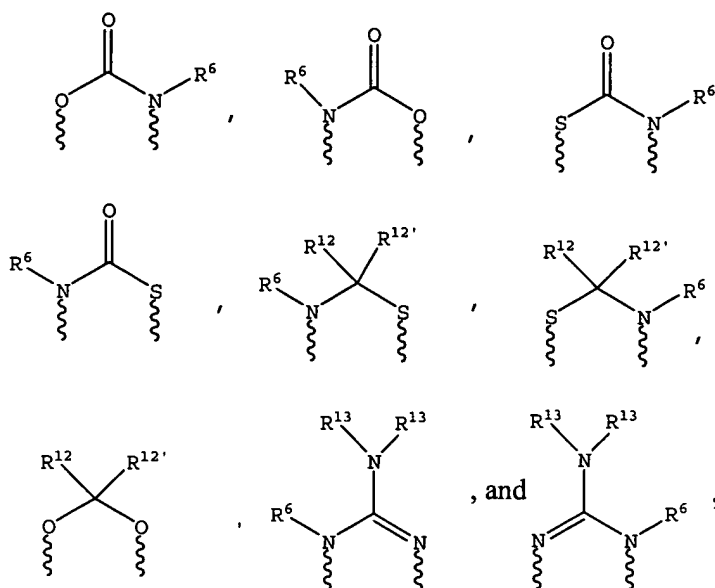
as to X, Y, and Z:

(a) one of X, Y, and Z is selected from the group consisting of C(O), S, S(O), S(O)₂, and NS(O)₂R⁷, and the remaining two of X, Y, and Z are CR⁸R⁹, and CR¹⁰R¹¹, or

(b) X and Z, or Z and Y together constitute a moiety selected from the group consisting of NR⁶C(O), NR⁶S(O), NR⁶S(O)₂, NR⁶S, NR⁶O, SS, NR⁶NR⁶, and OC(O), with the remaining one of X and Y being CR⁸R⁹, or

(c) n is zero and X, Y, and Z together constitute a moiety selected from the group consisting of:





wherein wavy lines are bonds to the atoms of the depicted ring;

R⁶ and R^{6'} are independently selected from the group consisting of hydrogen, formyl, sulfonic-C₁-C₆-alkyl, C₁-C₆-alkoxycarbonyl-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, C₁-C₆-alkylcarbonyl-C₁-C₆-alkyl, R⁸R⁹-aminocarbonyl-C₁-C₆-alkyl, C₁-C₆-alkoxycarbonyl-C₁-C₆-alkylcarbonyl, hydroxycarbonyl-C₁-C₆-alkylcarbonyl, C₁-C₆-alkylcarbonyl-C₁-C₆-alkylcarbonyl, C₁-C₆-alkoxycarbonylcarbonyl, hydroxycarbonylcarbonyl, C₁-C₆-alkylcarbonylcarbonyl, R⁸R⁹-aminocarbonylcarbonyl, C₁-C₆-alkanoyl, aryl-C₁-C₆-alkyl, aroyl, bis(C₁-C₆-alkoxy-C₁-C₆-alkyl)-C₁-C₆-alkyl, C₁-C₆-alkyl, C₁-C₆-haloalkyl, C₁-C₆-perfluoroalkyl, C₁-C₆-trifluoromethylalkyl, C₁-C₆-perfluoroalkoxy-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, C₃-C₆-cycloalkyl, heteroarylcarbonyl, heterocyclocarbonyl, 3- to 8-membered heterocycloalkyl, 3- to 8-membered heterocycloalkylcarbonyl, aryl, 5- to 6-membered heterocyclo, 5- to 6-membered heteroaryl, C₃-C₈-cycloalkyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, heteroaryl-C₁-C₆-alkoxy-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, arylsulfonyl, C₁-C₆-alkylsulfonyl, 5- to 6-membered heteroarylsulfonyl, carboxy-C₁-C₆-alkyl, C₁-C₄-alkoxycarbonyl-C₁-C₆-alkyl, aminocarbonyl, C₁-C₆-alkyl(R⁸N)iminocarbonyl,

aryl(R⁸N)iminocarbonyl, 5- to 6-membered heterocyclo(R⁸N)iminocarbonyl, arylthio-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, arylthio-C₃-C₆-alkenyl, C₁-C₄-alkylthio-C₃-C₆-alkenyl, 5- to 6-membered heteroaryl-C₁-C₆-alkyl, halo-C₁-C₆-alkanoyl, hydroxy-C₁-C₆-alkanoyl, thiol-C₁-C₆-alkanoyl, C₃-C₆-alkenyl, C₃-C₆-alkynyl, C₁-C₄-alkoxy-C₁-C₄-alkyl, C₁-C₅-alkoxycarbonyl, aryloxy carbonyl, NR⁸R⁹-(R⁸)iminomethyl, NR⁸R⁹-C₁-C₅-alkylcarbonyl, hydroxy-C₁-C₅-alkyl, R⁸R⁹-aminocarbonyl, R⁸R⁹-aminocarbonyl-C₁-C₆-alkylcarbonyl, hydroxyaminocarbonyl, R⁸R⁹-aminosulfonyl, R⁸R⁹-aminosulfon-C₁-C₆-alkyl, R⁸R⁹-amino-C₁-C₆-alkylsulfonyl, and R⁸R⁹-amino-C₁-C₆-alkyl;

R⁷ is selected from the group consisting of arylalkyl, aryl, heteroaryl, heterocyclo, C₁-C₆-alkyl, C₃-C₆-alkynyl, C₃-C₆-alkenyl, C₁-C₆-carboxyalkyl, and C₁-C₆-hydroxyalkyl; as to R⁸:

R⁸ is selected from the group consisting of hydrogen, hydroxy, C₁-C₆-alkyl, C₁-C₆-alkanoyl, aroyl, aryl, aryl-C₁-C₆-alkyl, heteroaryl, heteroaryl-C₁-C₆-alkyl, C₂-C₆-alkynyl, C₂-C₆-alkenyl, thiol-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, cycloalkyl, cycloalkyl-C₁-C₆-alkyl, heterocyclo-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, arylalkoxy-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkoxy-C₁-C₆-alkyl, hydroxy-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, hydroxycarbonylaryl-C₁-C₆-alkyl, aminocarbonyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, arylthio-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, C₁-C₆-alkylsulfinyl-C₁-C₆-alkyl, arylsulfinyl-C₁-C₆-alkyl, heteroarylsulfinyl-C₁-C₆-alkyl, C₁-C₆-alkylsulfonyl-C₁-C₆-alkyl, arylsulfonyl-C₁-C₆-alkyl, heteroarylsulfonyl-C₁-C₆-alkyl, perfluoro-C₁-C₆-alkyl, trifluoromethyl-C₁-C₆-alkyl, halo-C₁-C₆-alkyl, alkoxycarbonylamino-C₁-C₆-alkyl, and amino-C₁-C₆-alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C₁-C₆-alkyl, aryl-C₁-C₆-alkyl, cycloalkyl, and C₁-C₆-alkanoyl, or

R⁸ and R⁹, together with the carbon to which they are bonded, form a carbonyl group, or

R⁸ and R⁹ or R⁸ and R¹⁰, together with the atom(s) to which they are bonded, form a 5- to 8-membered carbocyclic ring, or a 5- to 8-membered heterocyclo or heteroaryl ring containing one or two heteroatoms independently selected from the group consisting of nitrogen, oxygen, and sulfur; as to R⁹:

R⁹ is selected from the group consisting of hydrogen, hydroxy, C₁-C₆-alkyl, C₁-C₆-alkanoyl, aroyl, aryl, aryl-C₁-C₆-alkyl, heteroaryl, heteroaryl-C₁-C₆-alkyl, C₂-C₆-alkynyl, C₂-C₆-alkenyl, thiol-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, cycloalkyl, cycloalkyl-C₁-C₆-alkyl, heterocyclo-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, arylalkoxy-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkoxy-C₁-C₆-alkyl, hydroxy-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, hydroxycarbonylaryl-C₁-C₆-alkyl, aminocarbonyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, arylthio-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, C₁-C₆-alkylsulfinyl-C₁-C₆-alkyl, arylsulfinyl-C₁-C₆-alkyl, heteroarylsulfinyl-C₁-C₆-alkyl, C₁-C₆-alkylsulfonyl-C₁-C₆-alkyl, arylsulfonyl-C₁-C₆-alkyl, heteroarylsulfonyl-C₁-C₆-alkyl, perfluoro-C₁-C₆-alkyl, trifluoromethyl-C₁-C₆-alkyl, halo-C₁-C₆-alkyl, alkoxycarbonylamino-C₁-C₆-alkyl, and amino-C₁-C₆-alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C₁-C₆-alkyl, aryl-C₁-C₆-alkyl, cycloalkyl, and C₁-C₆-alkanoyl, or

R^8 and R^9 , together with the carbon to which they are bonded, form a carbonyl group, a 5- to 8-membered carbocyclic ring, or a 5- to 8-membered heterocyclo or heteroaryl ring containing one or two heteroatoms independently selected from the group consisting of nitrogen, oxygen, and sulfur;
as to R^{10} :

R^{10} is selected from the group consisting of hydrogen, hydroxy, C_1 - C_6 -alkyl, C_1 - C_6 -alkanoyl, aroyl, aryl, aryl- C_1 - C_6 -alkyl, heteroaryl, heteroaryl- C_1 - C_6 -alkyl, C_2 - C_6 -alkynyl, C_2 - C_6 -alkenyl, thiol- C_1 - C_6 -alkyl, C_1 - C_6 -alkylthio- C_1 - C_6 -alkyl, cycloalkyl, cycloalkyl- C_1 - C_6 -alkyl, heterocyclo- C_1 - C_6 -alkyl, C_1 - C_6 -alkoxy- C_1 - C_6 -alkyl, arylalkoxy- C_1 - C_6 -alkyl, C_1 - C_6 -alkoxy- C_1 - C_6 -alkoxy- C_1 - C_6 -alkyl, hydroxy- C_1 - C_6 -alkyl, hydroxycarbonyl- C_1 - C_6 -alkyl, hydroxycarbonylaryl- C_1 - C_6 -alkyl, aminocarbonyl- C_1 - C_6 -alkyl, aryloxy- C_1 - C_6 -alkyl, heteroaryloxy- C_1 - C_6 -alkyl, arylthio- C_1 - C_6 -alkyl, heteroarylthio- C_1 - C_6 -alkyl, C_1 - C_6 -alkylsulfinyl- C_1 - C_6 -alkyl, arylsulfinyl- C_1 - C_6 -alkyl, heteroarylsulfinyl- C_1 - C_6 -alkyl, C_1 - C_6 -alkylsulfonyl- C_1 - C_6 -alkyl, arylsulfonyl- C_1 - C_6 -alkyl, heteroarylsulfonyl- C_1 - C_6 -alkyl, perfluoro- C_1 - C_6 -alkyl, trifluoromethyl- C_1 - C_6 -alkyl, halo- C_1 - C_6 -alkyl, alkoxycarbonylamino- C_1 - C_6 -alkyl, and amino- C_1 - C_6 -alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C_1 - C_6 -alkyl, aryl- C_1 - C_6 -alkyl, cycloalkyl, and C_1 - C_6 -alkanoyl,

R^{10} and R^{11} , together with the carbon to which they are bonded, form a carbonyl group, or

R^{10} and R^{11} or R^8 and R^{10} , together with the atom(s) to which they are bonded, form a 5- to 8-membered carbocyclic ring, or a 5- to 8-membered heterocyclo or heteroaryl ring containing one or two heteroatoms independently selected from the group consisting of nitrogen, oxygen, and sulfur;

as to R¹¹:

R¹¹ is selected from the group consisting of hydrogen, hydroxy, C₁-C₆-alkyl, C₁-C₆-alkanoyl, aroyl, aryl, aryl-C₁-C₆-alkyl, heteroaryl, heteroaryl-C₁-C₆-alkyl, C₂-C₆-alkynyl, C₂-C₆-alkenyl, thiol-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, cycloalkyl, cycloalkyl-C₁-C₆-alkyl, heterocyclo-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, arylalkoxy-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkoxy-C₁-C₆-alkyl, hydroxy-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, hydroxycarbonylaryl-C₁-C₆-alkyl, aminocarbonyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, arylthio-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, C₁-C₆-alkylsulfinyl-C₁-C₆-alkyl, arylsulfinyl-C₁-C₆-alkyl, heteroarylsulfinyl-C₁-C₆-alkyl, C₁-C₆-alkylsulfonyl-C₁-C₆-alkyl, arylsulfonyl-C₁-C₆-alkyl, heteroarylsulfonyl-C₁-C₆-alkyl, perfluoro-C₁-C₆-alkyl, trifluoromethyl-C₁-C₆-alkyl, halo-C₁-C₆-alkyl, alkoxycarbonylamino-C₁-C₆-alkyl, and amino-C₁-C₆-alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C₁-C₆-alkyl, aryl-C₁-C₆-alkyl, cycloalkyl, and C₁-C₆-alkanoyl, or

R¹⁰ and R¹¹, together with the carbon to which they are bonded, form a carbonyl group, a 5- to 8-membered carbocyclic ring, or a 5- to 8-membered heterocyclo or heteroaryl ring containing one or two heteroatoms independently selected from the group consisting of nitrogen, oxygen, and sulfur;

only one of R⁸ and R⁹ or R¹⁰ and R¹¹ is hydroxy;

R¹² and R^{12'} are independently selected from the group consisting of hydrogen, C₁-C₆-alkyl, aryl, aryl-C₁-C₆-alkyl, heteroaryl, heteroarylalkyl, C₂-C₆-alkynyl, C₂-C₆-alkenyl, thiol-C₁-C₆-alkyl, cycloalkyl, cycloalkyl-C₁-C₆-alkyl, heterocyclo-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, amino-C₁-C₆-alkyl,

C₁-C₆-alkoxy-C₁-C₆-alkoxy-C₁-C₆-alkyl, hydroxy-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, hydroxycarbonylaryl-C₁-C₆-alkyl, aminocarbonyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, arylthio-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, C₁-C₆-alkylsulfinyl-C₁-C₆-alkyl, arylsulfinyl-C₁-C₆-alkyl, heteroarylsulfinyl-C₁-C₆-alkyl, C₁-C₆-alkylsulfonyl-C₁-C₆-alkyl, arylsulfonyl-C₁-C₆-alkyl, heteroarylsulfonyl-C₁-C₆-alkyl, perfluoro-C₁-C₆-alkyl, trifluoromethyl-C₁-C₆-alkyl, halo-C₁-C₆-alkyl, alkoxycarbonylamino-C₁-C₆-alkyl, and amino-C₁-C₆-alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C₁-C₆-alkyl, aryl-C₁-C₆-alkyl, cycloalkyl, and C₁-C₆-alkanoyl;

R¹³ is selected from the group consisting of hydrogen, benzyl, phenyl, C₁-C₆-alkyl, C₂-C₆-alkynyl, C₂-C₆-alkenyl, and C₁-C₆-hydroxyalkyl;

Q is a 5- to 7-membered heterocyclic ring (other than piperazinyl) containing one or two nitrogen atoms;

A is selected from the group consisting of:

- (1) -O-,
- (2) -S-,
- (3) -NR¹⁷-,
- (4) -CO-N(R¹⁷),
- (5) -N(R¹⁷)-CO-,
- (6) -CO-O-,
- (7) -O-CO-,
- (8) -O-CO-O-,
- (9) -HC=CH-,
- (10) -NH-CO-NH-,
- (11) -C≡C-,

- (12) -NH-CO-O-,
- (13) -O-CO-NH-,
- (14) -N=N-,
- (15) -NH-NH-,
- (16) -CS-N(R¹⁷)-,
- (17) -N(R¹⁷)-CS-, and
- (18) a bond;

R¹⁷ is selected from the group consisting of hydrogen, C₁-C₄-alkyl, and phenyl;

R is selected from the group consisting of alkyl, alkoxyalkyl, aryl, heteroaryl, cycloalkyl, heterocyclo, arylalkyl, heteroarylalkyl, heterocycloalkyl, cycloalkylalkyl, cycloalkyloxyalkyl, heterocycloalkoxyalkyl, aryloxyalkyl, heteroaryloxyalkyl, arylthioalkyl, heteroarylthioalkyl, cycloalkylthioalkyl, and heterocycloalkylthioalkyl, wherein:

the aryl, heteroaryl, cycloalkyl, or heterocyclo is optionally substituted with up to two substituents independently selected from the group consisting of halo, alkyl, perfluoroalkyl, perfluoroalkoxy, perfluoroalkylthio, trifluoromethylalkyl, amino, alkoxycarbonylalkyl, alkoxy, C₁-C₂-alkylene-dioxy, hydroxycarbonylalkyl, hydroxycarbonylalkylamino, nitro, hydroxy, hydroxyalkyl, alkanoylamino, and alkoxycarbonyl;

R is other than alkyl or alkoxyalkyl when A is -O- or -S-;

E is selected from the group consisting of:

- (1) -CO(R¹⁹)-,
- (2) -(R¹⁹)CO-,
- (3) -CONH-,
- (4) -HNCO-,
- (5) -CO-,
- (6) -SO₂-R¹⁹-,
- (7) -R¹⁹-SO₂-,
- (8) -SO₂-,

- (9) -NH-SO₂-,
- (10) -SO₂-NH-,
- (11) -S-,
- (12) -NH-CO-O-,
- (13) -O-CO-NH-, and
- (14) a bond;

R¹⁹ is selected from the group consisting of heterocycloalkyl and cycloalkyl; and

Y² is selected from the group consisting of hydrogen, alkyl, alkoxy, haloalkyl, aryl, arylalkyl, cycloalkyl, heteroaryl, hydroxy, aryloxy, arylalkoxy, heteroaryloxy, heteroarylalkyl, perfluoroalkoxy, perfluoroalkylthio, trifluoromethylalkyl, alkenyl, heterocycloalkyl, trifluoromethyl, alkoxycarbonyl, and aminoalkyl, wherein:

the aryl, heteroaryl, aralkyl, or heterocycloalkyl is optionally substituted with up to two substituents independently selected from the group consisting of alkanoyl, halo, nitro, arylalkyl, aryl, alkoxy, trifluoroalkyl, trifluoroalkoxy, and amino, wherein:

the amino nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of alkyl and arylalkyl.

68. **(new)** A compound or salt according to claim 67, wherein g is 2.

69. **(new)** A compound or salt according to claim 68, wherein A-R-E-Y² is bonded at the 4-position of Q relative to the phenyl-bonded nitrogen of Q when Q is a 6- or 7-membered ring, and at the 3- or 4-position of Q relative to the phenyl-bonded nitrogen of Q when Q is a 5-membered ring.

70. **(new)** A compound or salt according to claim 69, wherein R¹⁴ is hydrogen.

71. **(new)** A compound or salt according to claim 69, wherein:

R¹⁴ is C(W)R²⁵;

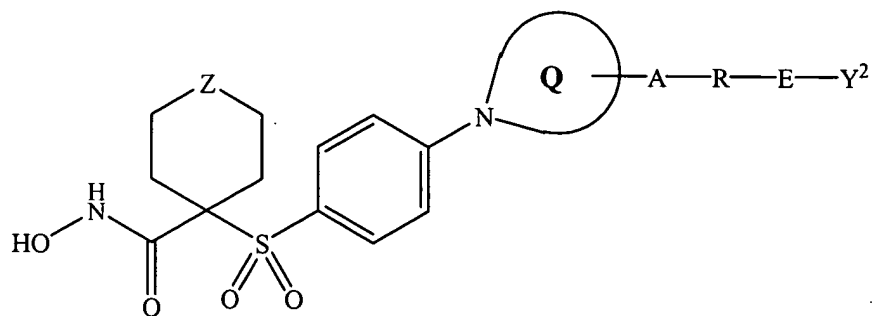
W is O; and

R^{25} is selected from the group consisting of C_1 - C_6 -alkyl, aryl, C_1 - C_6 -alkoxy, heteroaryl- C_1 - C_6 -alkyl, C_3 - C_8 -cycloalkyl- C_1 - C_6 -alkyl, and aryloxy.

72. **(new)** A compound or salt according to claim 69, wherein A is selected from the group consisting of -O- and -S-.

73. **(new)** A compound or salt according to claim 69, wherein R is selected from the group consisting of aryl, heteroaryl, cycloalkyl, and heterocyclo.

74. **(new)** A compound or salt according to claim 69, wherein the compound corresponds in structure to formula B-1A:

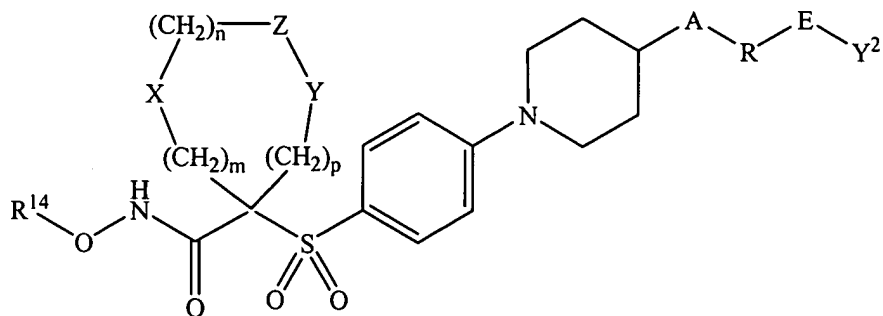


75. **(new)** A compound or salt according to claim 69, wherein Q is a 5-membered ring.

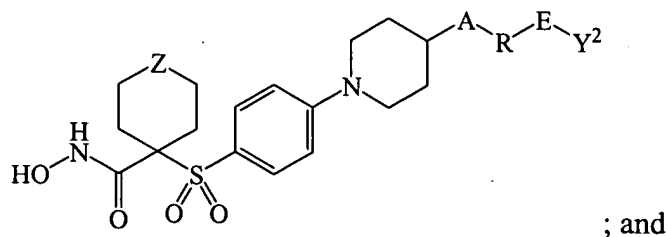
76. **(new)** A compound or salt according to claim 69, wherein Q is a 7-membered ring.

77. **(new)** A compound or salt according to claim 69, wherein Q is a 6-membered ring.

78. **(new)** A compound or salt according to claim 77, wherein the compound corresponds in structure to the following formula:



79. **(new)** A compound or salt according to claim 78, wherein:
 the compound corresponds in structure to the following formula:



Z is selected from the group consisting of C(O), S, S(O), S(O)₂, and NS(O)₂R⁷.

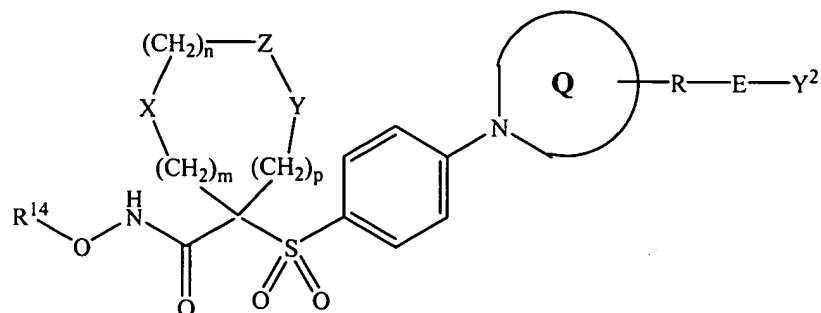
80. **(new)** A compound or salt according to claim 79, wherein R is selected from the group consisting of alkyl, alkoxyalkyl, aryl, heteroaryl, cycloalkyl, arylalkyl, heteroarylalkyl, heterocycloalkyl, cycloalkylalkyl, heterocycloalkoxyalkyl, aryloxyalkyl, heteroaryloxyalkyl, arylthioalkyl, heteroarylthioalkyl, and cycloalkylthioalkyl, wherein:

the aryl, heteroaryl, or cycloalkyl is optionally substituted with up to two substituents independently selected from the group consisting of halo, alkyl, perfluoroalkyl, perfluoroalkoxy, perfluoroalkylthio, trifluoromethylalkyl, amino, alkoxyalkyl, alkoxy, C₁-C₂-alkylene-dioxy, hydroxycarbonylalkyl, hydroxycarbonylalkylamino, nitro, hydroxy, hydroxyalkyl, alkanoylamino, and alkoxyalkyl.

81. **(new)** A compound or salt according to claim 80, wherein R is selected from the group consisting of aryl, heteroaryl, and cycloalkyl.

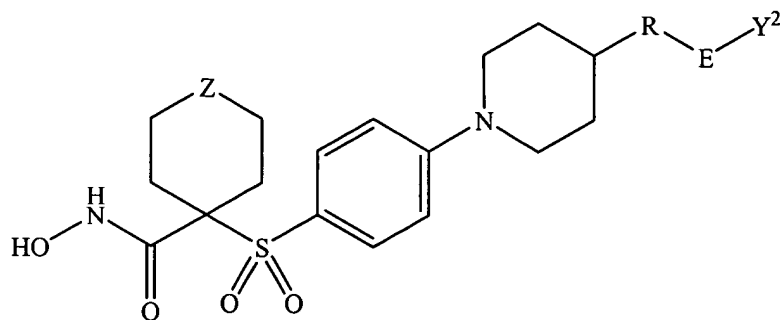
82. **(new)** A compound or salt according to claim 79, wherein Z is S.

83. **(new)** A compound or salt according to claim 68, wherein the compound corresponds in structure to the formula:



84. **(new)** A compound or salt according to claim 83, wherein R-E-Y² is bonded at the 4-position of Q relative to the phenyl-bonded nitrogen of Q when Q is a 6- or 7-membered ring, and at the 3- or 4-position of Q relative to the phenyl-bonded nitrogen of Q when Q is a 5-membered ring.

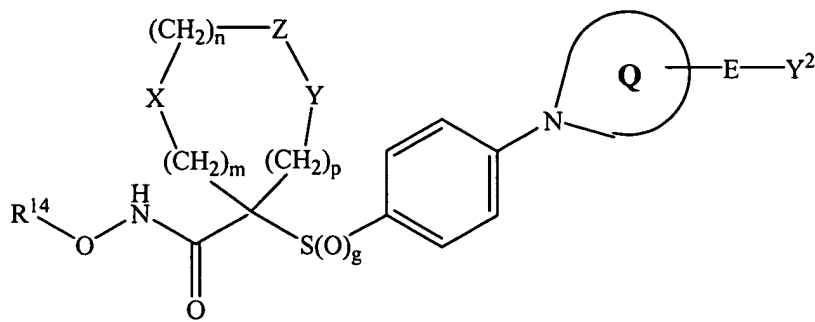
85. **(new)** A compound or salt according to claim 84, wherein:
the compound corresponds in structure to formula XI-1:



XI-1; and

Z is selected from the group consisting of C(O), S, S(O), S(O)₂, and NS(O)₂R⁷.

86. **(new)** A compound or a pharmaceutically acceptable salt thereof, wherein:
the compound corresponds in structure to the following formula:



R^{14} is selected from the group consisting of hydrogen and $C(W)R^{25}$;

W is selected from the group consisting of O and S;

R^{25} is selected from the group consisting of C_1 - C_6 -alkyl, aryl, C_1 - C_6 -alkoxy, heteroaryl- C_1 - C_6 -alkyl, C_3 - C_8 -cycloalkyl- C_1 - C_6 -alkyl, aryloxy, aryl- C_1 - C_6 -alkoxy, aryl- C_1 - C_6 -alkyl, heteroaryl, and amino C_1 - C_6 -alkyl, wherein:

the amino C_1 - C_6 -alkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C_1 - C_6 -alkyl, aryl, aryl- C_1 - C_6 -alkyl, C_3 - C_8 -cycloalkyl- C_1 - C_6 -alkyl, aryl- C_1 - C_6 -alkoxycarbonyl, C_1 - C_6 -alkoxycarbonyl, and C_1 - C_6 -alkanoyl, or

the amino C_1 - C_6 -alkyl nitrogen and two substituents attached thereto form a 5- to 8-membered heterocycle or heteroaryl ring;

g is zero, 1, or 2;

m is zero, 1, or 2;

n is zero, 1, or 2;

p is zero, 1, or 2;

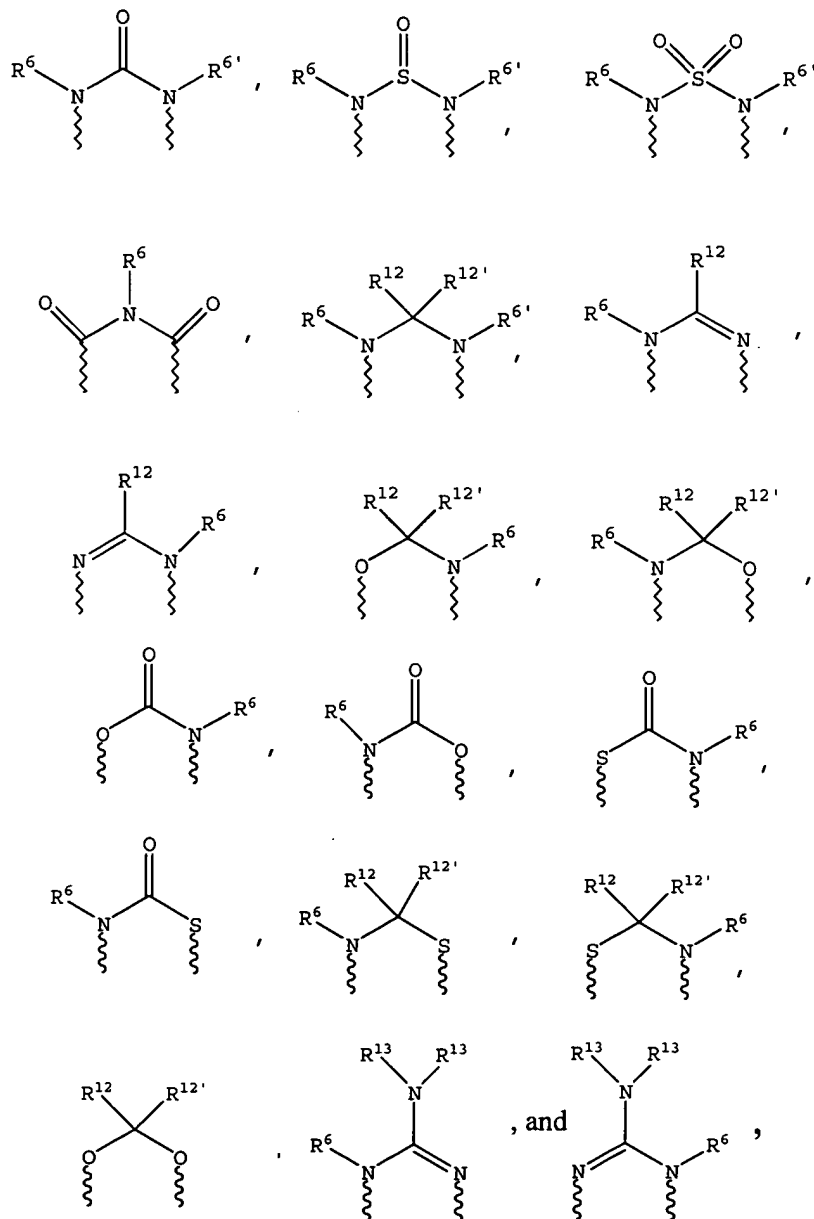
the sum of $m + n + p = 1, 2, 3$, or 4;

as to X, Y, and Z:

(a) one of X, Y, and Z is selected from the group consisting of C(O), S, S(O), S(O)₂, and NS(O)₂R⁷, and the remaining two of X, Y, and Z are CR⁸R⁹, and CR¹⁰R¹¹,
 or

(b) X and Z, or Z and Y together constitute a moiety selected from the group consisting of $\text{NR}^6\text{C}(\text{O})$, $\text{NR}^6\text{S}(\text{O})$, $\text{NR}^6\text{S}(\text{O})_2$, NR^6S , NR^6O , SS , NR^6NR^6 , and $\text{OC}(\text{O})$, with the remaining one of X and Y being CR^8R^9 , or

(c) n is zero and X, Y, and Z together constitute a moiety selected from the group consisting of:



wherein wavy lines are bonds to the atoms of the depicted ring;

R^6 and $\text{R}^{6'}$ are independently selected from the group consisting of hydrogen, formyl,

sulfonic-C₁-C₆-alkyl, C₁-C₆-alkoxycarbonyl-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, C₁-C₆-alkylcarbonyl-C₁-C₆-alkyl, R⁸R⁹-aminocarbonyl-C₁-C₆-alkyl, C₁-C₆-alkoxycarbonyl-C₁-C₆-alkylcarbonyl, hydroxycarbonyl-C₁-C₆-alkylcarbonyl, C₁-C₆-alkylcarbonyl-C₁-C₆-alkylcarbonyl, C₁-C₆-alkoxycarbonylcarbonyl, hydroxycarbonylcarbonyl, C₁-C₆-alkylcarbonylcarbonyl, R⁸R⁹-aminocarbonylcarbonyl, C₁-C₆-alkanoyl, aryl-C₁-C₆-alkyl, aroyl, bis(C₁-C₆-alkoxy-C₁-C₆-alkyl)-C₁-C₆-alkyl, C₁-C₆-alkyl, C₁-C₆-haloalkyl, C₁-C₆-perfluoroalkyl, C₁-C₆-trifluoromethylalkyl, C₁-C₆-perfluoroalkoxy-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, C₃-C₆-cycloalkyl, heteroarylcarbonyl, heterocyclocarbonyl, 3- to 8-membered heterocycloalkyl, 3- to 8-membered heterocycloalkylcarbonyl, aryl, 5- to 6-membered heterocyclo, 5- to 6-membered heteroaryl, C₃-C₈-cycloalkyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, heteroaryl-C₁-C₆-alkoxy-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, arylsulfonyl, C₁-C₆-alkylsulfonyl, 5- to 6-membered heteroarylsulfonyl, carboxy-C₁-C₆-alkyl, C₁-C₄-alkoxycarbonyl-C₁-C₆-alkyl, aminocarbonyl, C₁-C₆-alkyl(R⁸N)iminocarbonyl, aryl(R⁸N)iminocarbonyl, 5- to 6-membered heterocyclo(R⁸N)iminocarbonyl, arylthio-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, arylthio-C₃-C₆-alkenyl, C₁-C₄-alkylthio-C₃-C₆-alkenyl, 5- to 6-membered heteroaryl-C₁-C₆-alkyl, halo-C₁-C₆-alkanoyl, hydroxy-C₁-C₆-alkanoyl, thiol-C₁-C₆-alkanoyl, C₃-C₆-alkenyl, C₃-C₆-alkynyl, C₁-C₄-alkoxy-C₁-C₄-alkyl, C₁-C₅-alkoxycarbonyl, aryloxycarbonyl, NR⁸R⁹-(R⁸)iminomethyl, NR⁸R⁹-C₁-C₅-alkylcarbonyl, hydroxy-C₁-C₅-alkyl, R⁸R⁹-aminocarbonyl, R⁸R⁹-aminocarbonyl-C₁-C₆-alkylcarbonyl, hydroxyaminocarbonyl, R⁸R⁹-aminosulfonyl, R⁸R⁹-aminosulfon-C₁-C₆-alkyl, R⁸R⁹-amino-C₁-C₆-alkylsulfonyl, and R⁸R⁹-amino-C₁-C₆-alkyl;

R⁷ is selected from the group consisting of arylalkyl, aryl, heteroaryl, heterocyclo, C₁-C₆-alkyl, C₃-C₆-alkynyl, C₃-C₆-alkenyl, C₁-C₆-carboxyalkyl, and C₁-C₆-hydroxyalkyl;

as to R⁸:

R⁸ is selected from the group consisting of hydrogen, hydroxy, C₁-C₆-alkyl, C₁-C₆-alkanoyl, aroyl, aryl, aryl-C₁-C₆-alkyl, heteroaryl, heteroaryl-C₁-C₆-alkyl, C₂-C₆-alkynyl, C₂-C₆-alkenyl, thiol-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, cycloalkyl, cycloalkyl-C₁-C₆-alkyl, heterocyclo-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, arylalkoxy-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkoxy-C₁-C₆-alkyl, hydroxy-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, hydroxycarbonylaryl-C₁-C₆-alkyl, aminocarbonyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, arylthio-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, C₁-C₆-alkylsulfinyl-C₁-C₆-alkyl, arylsulfinyl-C₁-C₆-alkyl, heteroarylsulfinyl-C₁-C₆-alkyl, C₁-C₆-alkylsulfonyl-C₁-C₆-alkyl, arylsulfonyl-C₁-C₆-alkyl, heteroarylsulfonyl-C₁-C₆-alkyl, perfluoro-C₁-C₆-alkyl, trifluoromethyl-C₁-C₆-alkyl, halo-C₁-C₆-alkyl, alkoxycarbonylamino-C₁-C₆-alkyl, and amino-C₁-C₆-alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C₁-C₆-alkyl, aryl-C₁-C₆-alkyl, cycloalkyl, and C₁-C₆-alkanoyl, or

R⁸ and R⁹, together with the carbon to which they are bonded, form a carbonyl group, or

R⁸ and R⁹ or R⁸ and R¹⁰, together with the atom(s) to which they are bonded, form a 5- to 8-membered carbocyclic ring, or a 5- to 8-membered heterocyclo or heteroaryl ring containing one or two heteroatoms independently selected from the group consisting of nitrogen, oxygen, and sulfur;

as to R⁹:

R⁹ is selected from the group consisting of hydrogen, hydroxy, C₁-C₆-alkyl, C₁-C₆-alkanoyl, aroyl, aryl, aryl-C₁-C₆-alkyl, heteroaryl, heteroaryl-C₁-C₆-alkyl,

C₂-C₆-alkynyl, C₂-C₆-alkenyl, thiol-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, cycloalkyl, cycloalkyl-C₁-C₆-alkyl, heterocyclo-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, arylalkoxy-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkoxy-C₁-C₆-alkyl, hydroxy-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, hydroxycarbonylaryl-C₁-C₆-alkyl, aminocarbonyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, arylthio-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, C₁-C₆-alkylsulfinyl-C₁-C₆-alkyl, arylsulfinyl-C₁-C₆-alkyl, heteroarylsulfinyl-C₁-C₆-alkyl, C₁-C₆-alkylsulfonyl-C₁-C₆-alkyl, arylsulfonyl-C₁-C₆-alkyl, heteroarylsulfonyl-C₁-C₆-alkyl, perfluoro-C₁-C₆-alkyl, trifluoromethyl-C₁-C₆-alkyl, halo-C₁-C₆-alkyl, alkoxycarbonylamino-C₁-C₆-alkyl, and amino-C₁-C₆-alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C₁-C₆-alkyl, aryl-C₁-C₆-alkyl, cycloalkyl, and C₁-C₆-alkanoyl, or

R⁸ and R⁹, together with the carbon to which they are bonded, form a carbonyl group, a 5- to 8-membered carbocyclic ring, or a 5- to 8-membered heterocyclo or heteroaryl ring containing one or two heteroatoms independently selected from the group consisting of nitrogen, oxygen, and sulfur; as to R¹⁰:

R¹⁰ is selected from the group consisting of hydrogen, hydroxy, C₁-C₆-alkyl, C₁-C₆-alkanoyl, aroyl, aryl, aryl-C₁-C₆-alkyl, heteroaryl, heteroaryl-C₁-C₆-alkyl, C₂-C₆-alkynyl, C₂-C₆-alkenyl, thiol-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, cycloalkyl, cycloalkyl-C₁-C₆-alkyl, heterocyclo-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, arylalkoxy-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkoxy-C₁-C₆-alkyl, hydroxy-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, hydroxycarbonylaryl-C₁-C₆-alkyl,

aminocarbonyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, arylthio-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, C₁-C₆-alkylsulfinyl-C₁-C₆-alkyl, arylsulfinyl-C₁-C₆-alkyl, heteroarylsulfinyl-C₁-C₆-alkyl, C₁-C₆-alkylsulfonyl-C₁-C₆-alkyl, arylsulfonyl-C₁-C₆-alkyl, heteroarylsulfonyl-C₁-C₆-alkyl, perfluoro-C₁-C₆-alkyl, trifluoromethyl-C₁-C₆-alkyl, halo-C₁-C₆-alkyl, alkoxy-carbonylamino-C₁-C₆-alkyl, and amino-C₁-C₆-alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of C₁-C₆-alkyl, aryl-C₁-C₆-alkyl, cycloalkyl, and C₁-C₆-alkanoyl,

R¹⁰ and R¹¹, together with the carbon to which they are bonded, form a carbonyl group, or

R¹⁰ and R¹¹ or R⁸ and R¹⁰, together with the atom(s) to which they are bonded, form a 5- to 8-membered carbocyclic ring, or a 5- to 8-membered heterocyclo or heteroaryl ring containing one or two heteroatoms independently selected from the group consisting of nitrogen, oxygen, and sulfur; as to R¹¹:

R¹¹ is selected from the group consisting of hydrogen, hydroxy, C₁-C₆-alkyl, C₁-C₆-alkanoyl, aroyl, aryl, aryl-C₁-C₆-alkyl, heteroaryl, heteroaryl-C₁-C₆-alkyl, C₂-C₆-alkynyl, C₂-C₆-alkenyl, thiol-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl, cycloalkyl, cycloalkyl-C₁-C₆-alkyl, heterocyclo-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, arylalkoxy-C₁-C₆-alkyl, C₁-C₆-alkoxy-C₁-C₆-alkoxy-C₁-C₆-alkyl, hydroxy-C₁-C₆-alkyl, hydroxycarbonyl-C₁-C₆-alkyl, hydroxycarbonylaryl-C₁-C₆-alkyl, aminocarbonyl-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, arylthio-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, C₁-C₆-alkylsulfinyl-C₁-C₆-alkyl, arylsulfinyl-C₁-C₆-alkyl, heteroarylsulfinyl-C₁-C₆-alkyl,

C₁-C₆-alkylsulfonyl-C₁-C₆-alkyl, arylsulfonyl-C₁-C₆-alkyl,
heteroarylsulfonyl-C₁-C₆-alkyl, perfluoro-C₁-C₆-alkyl, trifluoromethyl-C₁-C₆-alkyl,
halo-C₁-C₆-alkyl, alkoxycarbonylamino-C₁-C₆-alkyl, and amino-C₁-C₆-alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two
substituents independently selected from the group consisting of C₁-C₆-alkyl,
aryl-C₁-C₆-alkyl, cycloalkyl, and C₁-C₆-alkanoyl, or

R¹⁰ and R¹¹, together with the carbon to which they are bonded, form a carbonyl
group, a 5- to 8-membered carbocyclic ring, or a 5- to 8-membered heterocyclo or
heteroaryl ring containing one or two heteroatoms independently selected from the group
consisting of nitrogen, oxygen, and sulfur;

only one of R⁸ and R⁹ or R¹⁰ and R¹¹ is hydroxy;

R¹² and R^{12'} are independently selected from the group consisting of hydrogen,
C₁-C₆-alkyl, aryl, aryl-C₁-C₆-alkyl, heteroaryl, heteroarylalkyl, C₂-C₆-alkynyl, C₂-C₆-alkenyl,
thiol-C₁-C₆-alkyl, cycloalkyl, cycloalkyl-C₁-C₆-alkyl, heterocyclo-C₁-C₆-alkyl,
C₁-C₆-alkoxy-C₁-C₆-alkyl, aryloxy-C₁-C₆-alkyl, amino-C₁-C₆-alkyl,
C₁-C₆-alkoxy-C₁-C₆-alkoxy-C₁-C₆-alkyl, hydroxy-C₁-C₆-alkyl,
hydroxycarbonyl-C₁-C₆-alkyl, hydroxycarbonylaryl-C₁-C₆-alkyl, aminocarbonyl-C₁-C₆-alkyl,
aryloxy-C₁-C₆-alkyl, heteroaryloxy-C₁-C₆-alkyl, C₁-C₆-alkylthio-C₁-C₆-alkyl,
arylthio-C₁-C₆-alkyl, heteroarylthio-C₁-C₆-alkyl, C₁-C₆-alkylsulfinyl-C₁-C₆-alkyl,
arylsulfinyl-C₁-C₆-alkyl, heteroarylsulfinyl-C₁-C₆-alkyl, C₁-C₆-alkylsulfonyl-C₁-C₆-alkyl,
arylsulfonyl-C₁-C₆-alkyl, heteroarylsulfonyl-C₁-C₆-alkyl, perfluoro-C₁-C₆-alkyl,
trifluoromethyl-C₁-C₆-alkyl, halo-C₁-C₆-alkyl, alkoxycarbonylamino-C₁-C₆-alkyl, and
amino-C₁-C₆-alkyl, wherein:

the aminoalkyl nitrogen is optionally substituted with up to two substituents
independently selected from the group consisting of C₁-C₆-alkyl, aryl-C₁-C₆-alkyl,
cycloalkyl, and C₁-C₆-alkanoyl;

R^{13} is selected from the group consisting of hydrogen, benzyl, phenyl, C_1 - C_6 -alkyl, C_2 - C_6 -alkynyl, C_2 - C_6 -alkenyl, and C_1 - C_6 -hydroxyalkyl;

Q is a 5- to 7-membered heterocyclic ring (other than piperazinyl) containing one or two nitrogen atoms;

E is selected from the group consisting of:

- (1) $-\text{CO}(R^{19})-$,
- (2) $-(R^{19})\text{CO}-$,
- (3) $-\text{CONH}-$,
- (4) $-\text{HNCO}-$,
- (5) $-\text{CO}-$,
- (6) $-\text{SO}_2\text{-}R^{19}-$,
- (7) $\text{-}R^{19}\text{-SO}_2-$,
- (8) $-\text{SO}_2-$,
- (9) $-\text{NH-SO}_2-$,
- (10) $-\text{SO}_2\text{-NH}-$,
- (11) $-\text{S}-$,
- (12) $-\text{NH-CO-O}-$,
- (13) $-\text{O-CO-NH}-$, and
- (14) a bond;

R^{19} is selected from the group consisting of heterocycloalkyl and cycloalkyl; and

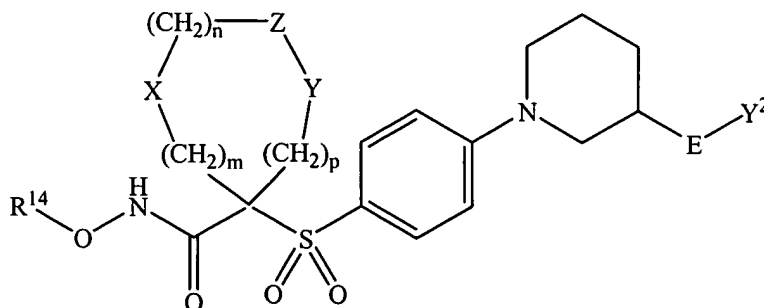
Y^2 is selected from the group consisting of alkyl, alkoxy, haloalkyl, aryl, arylalkyl, cycloalkyl, heteroaryl, hydroxy, aryloxy, arylalkoxy, heteroaryloxy, heteroarylalkyl, perfluoroalkoxy, perfluoroalkylthio, trifluoromethylalkyl, alkenyl, heterocyclo, heterocycloalkyl, trifluoromethyl, alkoxycarbonyl, and aminoalkyl, wherein:

the aryl, heteroaryl, aralkyl, heterocyclo, or heterocycloalkyl is optionally substituted with up to two substituents independently selected from the group consisting of alkyl, alkanoyl, halo, nitro, arylalkyl, aryl, alkoxy, trifluoroalkyl, trifluoroalkoxy, and amino, wherein:

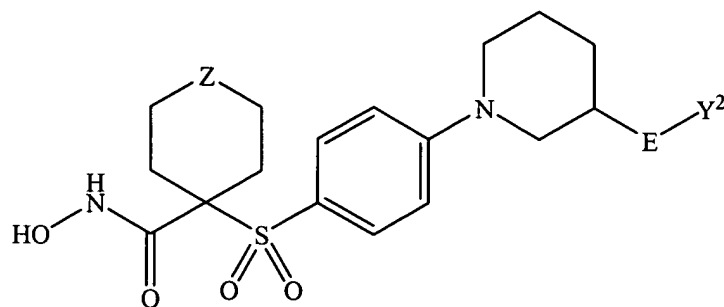
the amino nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of alkyl and arylalkyl.

87. **(new)** A compound or salt according to claim 86, wherein g is 2.

88. **(new)** A compound or salt according to claim 87, wherein the compound corresponds in structure to the following formula:



89. **(new)** A compound or salt according to claim 87, wherein:
 the compound corresponds in structure to formula IX-2:



IX-2; and

Z is selected from the group consisting of C(O), S, S(O), S(O)₂, and NS(O)₂R⁷.

90. **(new)** A compound or salt according to claim 87, wherein E-Y² is bonded at the 4-position of Q relative to the phenyl-bonded nitrogen of Q when Q is a 6- or 7-membered ring, and at the 3- or 4-position of Q relative to the phenyl-bonded nitrogen of Q when Q is a 5-membered ring.

91. (new) A compound or salt according to claim 90, wherein Y^2 is selected from the group consisting of alkyl, alkoxy, haloalkyl, aryl, arylalkyl, cycloalkyl, heteroaryl, hydroxy, aryloxy, arylalkoxy, heteroaryloxy, heteroarylalkyl, perfluoroalkoxy, perfluoroalkylthio, trifluoromethylalkyl, alkenyl, heterocycloalkyl, trifluoromethyl, alkoxycarbonyl, and aminoalkyl, wherein:

the aryl, heteroaryl, aralkyl, or heterocycloalkyl is optionally substituted with up to two substituents independently selected from the group consisting of alkanoyl, halo, nitro, arylalkyl, aryl, alkoxy, trifluoroalkyl, trifluoroalkoxy, and amino, wherein:

the amino nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of alkyl and arylalkyl.

92. (new) A compound or salt according to claim 90, wherein R^{14} is hydrogen.

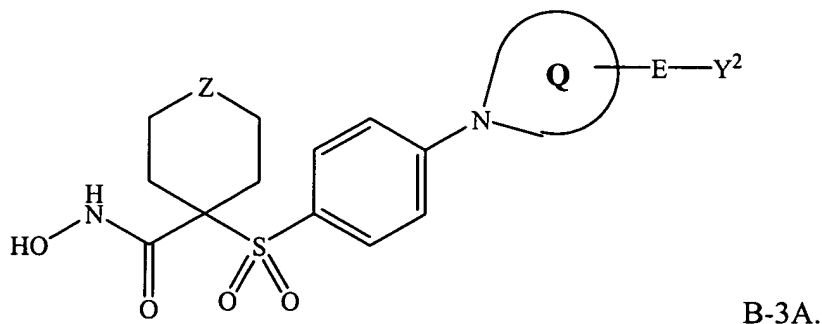
93. (new) A compound or salt according to claim 90, wherein:

R^{14} is $C(W)R^{25}$;

W is O; and

R^{25} is selected from the group consisting of C_1 - C_6 -alkyl, aryl, C_1 - C_6 -alkoxy, heteroaryl- C_1 - C_6 -alkyl, C_3 - C_8 -cycloalkyl- C_1 - C_6 -alkyl, and aryloxy.

94. (new) A compound or salt according to claim 90, wherein the compound corresponds in structure to formula B-3A:

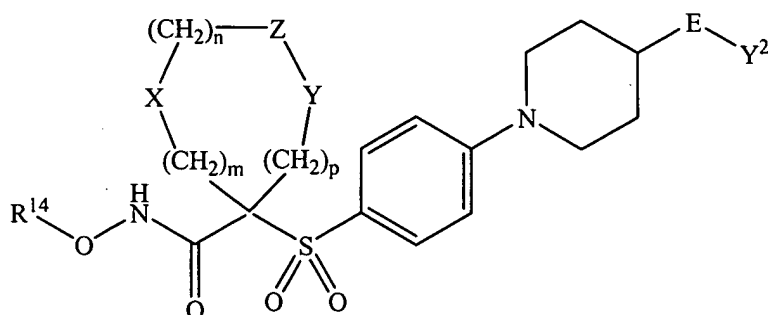


95. (new) A compound or salt according to claim 90, wherein Q is a 5-membered ring.

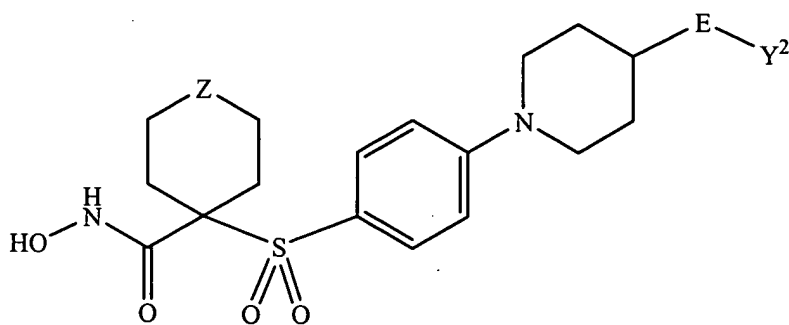
96. (new) A compound or salt according to claim 90, wherein Q is a 7-membered ring.

97. (new) A compound or salt according to claim 90, wherein Q is a 6-membered ring.

98. (new) A compound or salt according to claim 97, wherein the compound corresponds in structure to the following formula:



99. (new) A compound or salt according to claim 98, wherein:
the compound corresponds in structure to formula IX-1:



IX-1; and

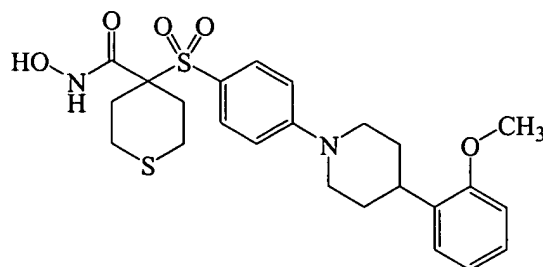
Z is selected from the group consisting of C(O), S, S(O), S(O)₂, and NS(O)₂R⁷.

100. (new) A compound or salt according to claim 99, wherein Y² is selected from the group consisting of alkyl, alkoxy, haloalkyl, aryl, arylalkyl, cycloalkyl, heteroaryl, hydroxy, aryloxy, arylalkoxy, heteroaryloxy, heteroarylalkyl, perfluoroalkoxy, perfluoroalkylthio, trifluoromethylalkyl, alkenyl, heterocycloalkyl, trifluoromethyl, alkoxycarbonyl, and aminoalkyl, wherein:

the aryl, heteroaryl, aralkyl, or heterocycloalkyl is optionally substituted with up to two substituents independently selected from the group consisting of alkanoyl, halo, nitro, arylalkyl, aryl, alkoxy, trifluoroalkyl, trifluoroalkoxy, and amino, wherein: the amino nitrogen is optionally substituted with up to two substituents independently selected from the group consisting of alkyl and arylalkyl.

101. **(new)** A compound or salt according to claim 99, wherein Z is S.

102. **(new)** A compound or salt according to claim 101, wherein the compound corresponds in structure to the formula:



103. **(new)** A method for treating a pathological condition in a mammal, wherein: the condition is treatable by inhibiting matrix metalloprotease activity; the method comprises administering a compound or a pharmaceutically acceptable salt thereof to the mammal in an amount that is effective to treat the condition; the compound corresponds in structure to a compound recited in claim 68; and the compound or salt inhibits the activity of one or more of MMP-2, MMP-9, and MMP-13, while exhibiting substantially less inhibitory activity against MMP-1.

104. **(new)** A method according to claim 103, wherein A is a bond.

105. **(new)** A method according to claim 103, wherein the pathological condition is selected from the group consisting of rheumatoid arthritis, osteoarthritis, septic arthritis, corneal, epidermal or gastric ulceration, tumor metastasis, invasion or angiogenesis, periodontal disease, proteinuria, Alzheimer's disease, coronary thrombosis, and bone disease.

106. **(new)** A method according to claim 105, wherein the pathological condition is osteoarthritis.

107. **(new)** A method according to claim 103, wherein the compound or salt is administered a plurality of times.

108. **(new)** A method for treating a pathological condition in a mammal, wherein:
the condition is treatable by inhibiting matrix metalloprotease activity;
the method comprises administering a compound or a pharmaceutically acceptable salt thereof to the mammal in an amount that is effective to treat the condition;
the compound corresponds in structure to a compound recited in claim 87; and
the compound or salt inhibits the activity of one or more of MMP-2, MMP-9, and MMP-13, while exhibiting substantially less inhibitory activity against MMP-1.

109. **(new)** A method according to claim 108, wherein the pathological condition is selected from the group consisting of rheumatoid arthritis, osteoarthritis, septic arthritis, corneal, epidermal or gastric ulceration, tumor metastasis, invasion or angiogenesis, periodontal disease, proteinuria, Alzheimer's disease, coronary thrombosis, and bone disease.

110. **(new)** A method according to claim 109, wherein the pathological condition is osteoarthritis.

111. **(new)** A method according to claim 108, wherein the compound or salt is administered a plurality of times.

112. **(new)** A pharmaceutical composition that comprises a compound or salt according to claim 68 dissolved or dispersed in a pharmaceutically acceptable carrier.

113. **(new)** A pharmaceutical composition according to claim 112, wherein A is a bond.

114. (new) A pharmaceutical composition that comprises a compound or salt according to claim 87 dissolved or dispersed in a pharmaceutically acceptable carrier.